

AUTOMATED FINITE ELEMENT GRID BREAK-UP METHOD - A

VERIFICATION OF THE SIX NODE AVERAGING APPROACH

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SUMMARY

The verification of the nodal averaging method for generating automated finite triangular element grids has been demonstrated. This was accomplished with a six node averaging program (SNAP) which was placed on an IBM 2250 vector graphics scope terminal. The advantage of this method is that it is unnecessary to program time consuming geometric division and transition algorithms.

INTRODUCTION

The most popular method for the automatic break-up of regions into plane finite elements seems to be variations of the "sides-and-parts" method as discussed in references 1 through 4. In this method the structure is divided into four-sided parts with the sides divided into straight line segments connected at nodes. Figures 1, 2, and 3 show such parts in the form of a square which has a one-to-one topological correspondence to a general four-sided figure. If the number of nodes on pairs of opposite sides are equal, lines or curves drawn through corresponding side nodes will generate internal mesh points at their intersections, as shown in figure 1.

For further refinement a division algorithm can be used to specify variable spacing in both directions. If the nodes on one pair of sides are equal in number, while those on the other pair of sides are unequal, this algorithm can be used to generate a transition pattern as shown in figure 2. In this instance the algorithm is that the number of nodes on a row is one less or one more than the number on an adjacent row. In both of these cases the topological pattern is easily predictable and is amenable to automatic mesh generation. If the numbers of nodes on both pairs of sides are not equal, the pattern is generally not too apparent, although exceptions do occur, as shown in figures 3 and 4. The pattern in figure 3 is obtained by drawing the full lines and completing the triangulation by connecting the mesh points at their intersections, as shown by the dashed lines. Here there is a transition from four to five nodes on one pair of opposite sides and a transition from three to four nodes on the other pair. This pattern could have been achieved by using the transition algorithm in two directions and a little imagination. Figure 4 shows the pattern for the same configuration which was obtained by the method which is described herein. This pattern was obtained without specifying

division or transition algorithms and depends only upon the positions of the side nodes. Note that it has the same topological pattern as that shown in figure 3; i. e., the nodes in both figures can be made coincident by shifting without destroying their connectivity.

A more general case in which the division or transition algorithms are more difficult to formulate is shown in figure 5. This pattern was obtained again by the method which is described herein and again required the definition of the position of the side nodes only. Here, the division and transition algorithms are automatic. By inspection it can be seen that the number of nodes along any interior segmented curve is one more or less than the number on adjacent curves, and hence the transition algorithm is satisfied. The transition pattern for this configuration could also have been predicted with some imagination and hind-sight and computerized. However, as the number of side nodes increases, the transition patterns become more difficult to formulate for computer programming.

NODE AVERAGING APPROACH

This technique was programmed because it makes it unnecessary to use transition and division algorithms or to input internal paths or subsystems. The only rigid condition that is stipulated when using node averaging is that each internal mesh point always be the common vertex of six triangular elements. This is stipulated in order to simplify the computer formulation and to provide uniform nodal patterns at each internal mesh point. If this requirement is met, the topological structure of the break-up is a series of nested polygons each of which has a multiple of six nodes or edges whose number differs by six from the number in adjacent polygons. The reason that this pattern generates the desired transition algorithm is quite obvious if one examines the sector ABCDEF in figure 6. The number of nodes across the sector increases in increments of 1 as the boundaries of the nested polygons are crossed. Thus, there is one node at A, two nodes along BE, three nodes along CF and finally, four nodes along DG. This algorithm exists in each of the six sectors. The triangulation is achieved by connecting nodes in adjacent polygons as shown in figure 6. Since the external contour of the region is the extreme polygon, the number of known contour nodes or edges also must always be specified as a multiple of six.

We now treat the coordinates of the interior nodes as unknowns and stipulate that they are the average of the coordinates of the six adjacent nodes. This is a rough application of the finite difference method for solving Laplace's equation for a region with prescribed boundary conditions. This algorithm yields two sets of simultaneous equations in which the coordinates of the internal nodes are the unknowns. If the number of nodes is too large for simultaneous equation solutions, the coordinates can be obtained by assuming initial values and iterating until convergence is attained.

The program which has been developed uses the simultaneous equation method. If a hole is present in the region, the number of nodes on the hole contour must also be a multiple of six. The contour of the hole then becomes the first nested polygon while the external contour becomes the last polygon or vice versa. Since the number of nodes on each polygon follows an arithmetic series, the total number of nodes and elements can be expressed in terms of the number

of internal and external contour nodes. If N_e is the number of external nodes and N_i is the number of internal nodes, these relations are:

$$\text{Total number of nodes} = \frac{N_e + N_i}{2} + \frac{N_e^2 - N_i^2}{12}$$

$$\text{Total number of elements} = \frac{N_e^2 - N_i^2}{6}$$

If no hole is present, the relations are:

$$\text{Total number of nodes} = 1 + \frac{N_e}{2} + \frac{N_e^2}{12}$$

$$\text{Total number of elements} = \frac{N_e^2}{6}$$

Figures 4 and 5 show the patterns obtained by solving the set of linear simultaneous equations using the node averaging algorithm. Attempts to use this algorithm on regions with concave and/or slender contours were generally unsuccessful because mesh points had a tendency to be near a contour or outside the region. Such regions can be handled by dividing them into subregions and piecing together independent solutions. This technique, however, negates the advantage of a single break-up. Other methods of handling concave boundaries are discussed in the references.

This automatic break-up algorithm has been programmed into a finite element structures program and a compatible finite element break-up scope program. This requires the input of contour nodes only as discussed previously. A test case was input and run on the scope program and a plot of the sequence of presentations is shown in figures 7 through 11. The first display (figure 7) shows the contour nodes generated by inputting corner nodes, hole location, and hole diameter while using an automatic boundary break-up option, which is also incorporated in the program. The second display (figure 8) is the generated automatic break-up. The third display (figure 9) is a refined break-up which is obtained by triangulating the mid-points of the edges of each element in the previous coarse break-up. The fourth display (figure 10) shows the coarse break-up as modified by moving nodes with the light-pen. The fifth display (figure 11) is the fine break-up obtained by triangulating the mid-points of the edges in the previous adjusted break-up. The light-pen can also be used to present magnified local regions, sub-systems, and boundary and perimeter nodes.

CONCLUDING REMARKS

A nodal pattern for triangular mesh generation, using nodal averaging techniques, has been developed. This pattern has intrinsic properties which make it unnecessary to program internal transition algorithms. An automated mesh generation program using this technique has been developed for use on the

IBM 2250 vector graphics scope terminal. The use of this program on this terminal allows the operator to modify geometry in order to improve the break-up prior to being input into an associated finite element structures program.

REFERENCES

1. Buell, W. R., and B. A. Bush: Mesh Generation - A Survey. *Transactions of ASME*, February 1973, pp. 332-338.
2. Crose, J. G., and R. M. Jones: SAAS II, Finite Element Stress Analysis of Axisymmetric Solids With Orthotropic, Temperature-Dependent Material Properties. TR-0200 (S4980)-1, The Aerospace Corporation, September 1968.
3. Frederick, C. O., Y. C. Wong, and F. W. Edge: Two-Dimensional Automatic Mesh Generation for Structural Analysis. *International Journal for Numerical Methods in Engineering*, 1970, Volume 2, pp. 133-144.
4. Jones, R. E.: QMESH: A Self-Organizing Mesh Generation Program. SLA-73-1088, Sandia Laboratories, July 1974.

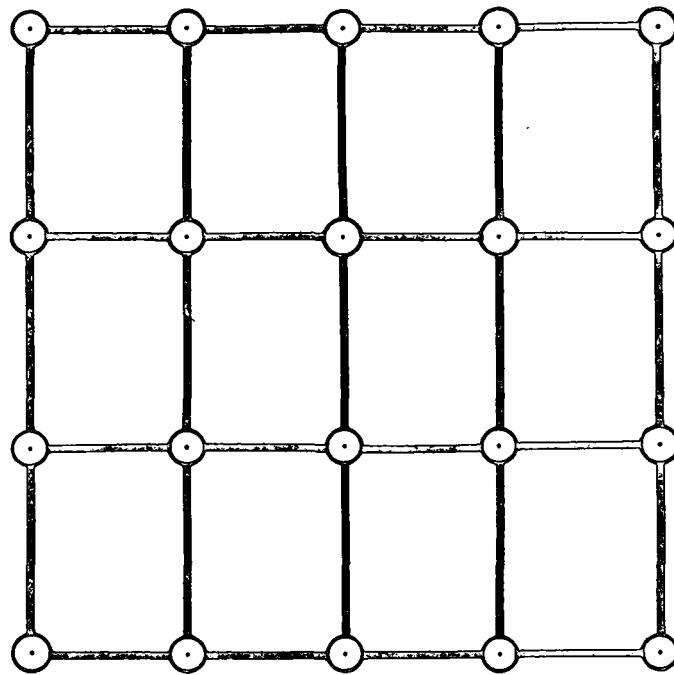


Figure 1. Conventional simple break-up.

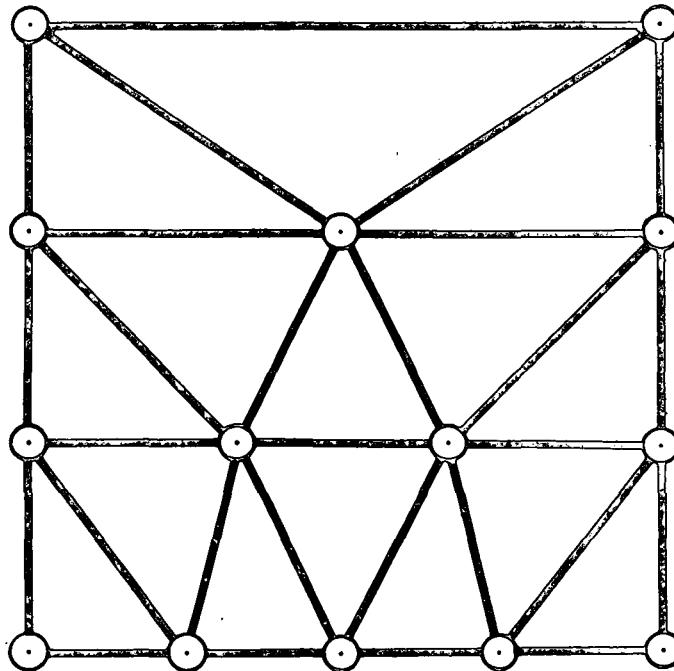


Figure 2. Conventional break-up with single transition.

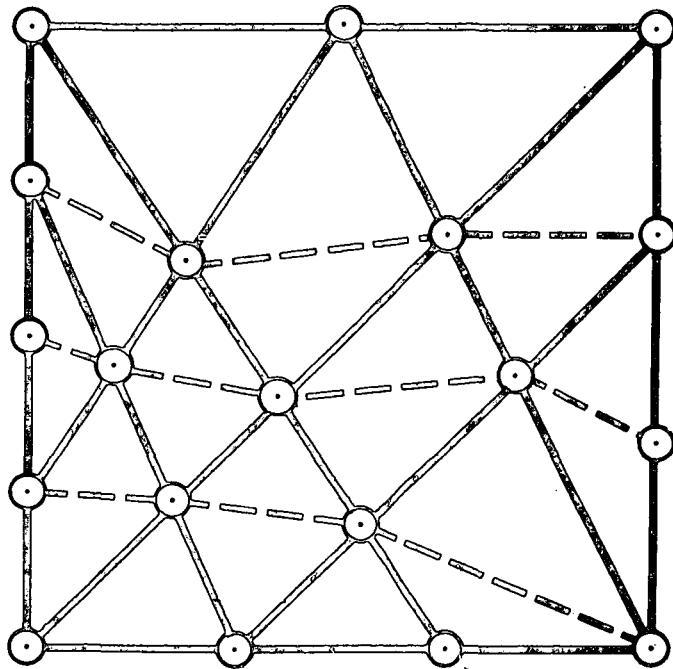


Figure 3. Double transition break-up.

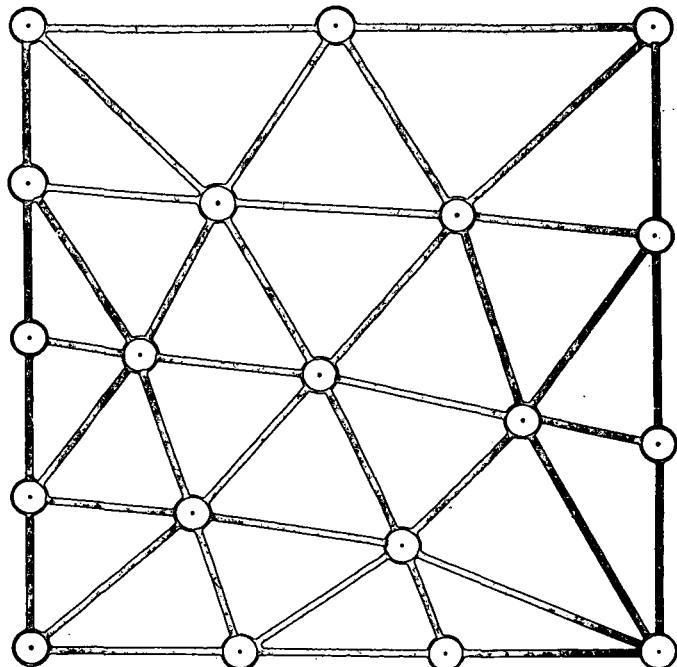


Figure 4. SNAP break-up.

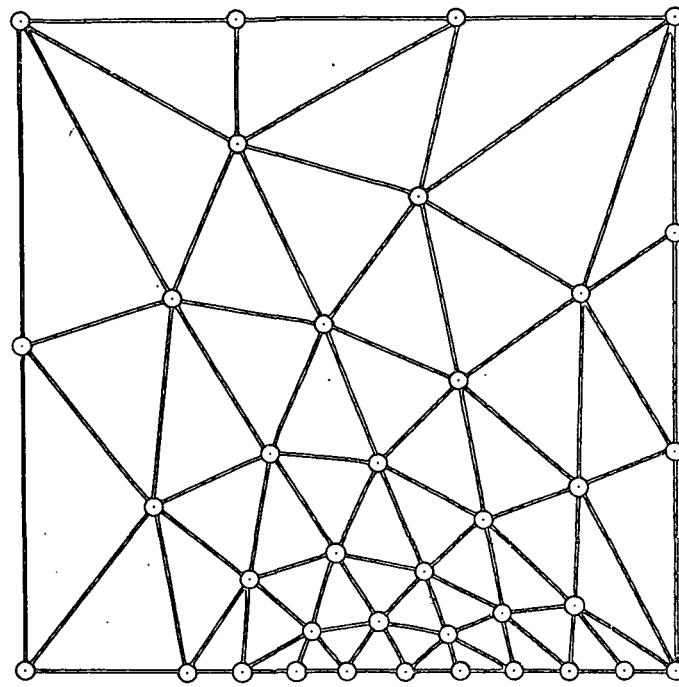


Figure 5. Complex SNAP break-up.

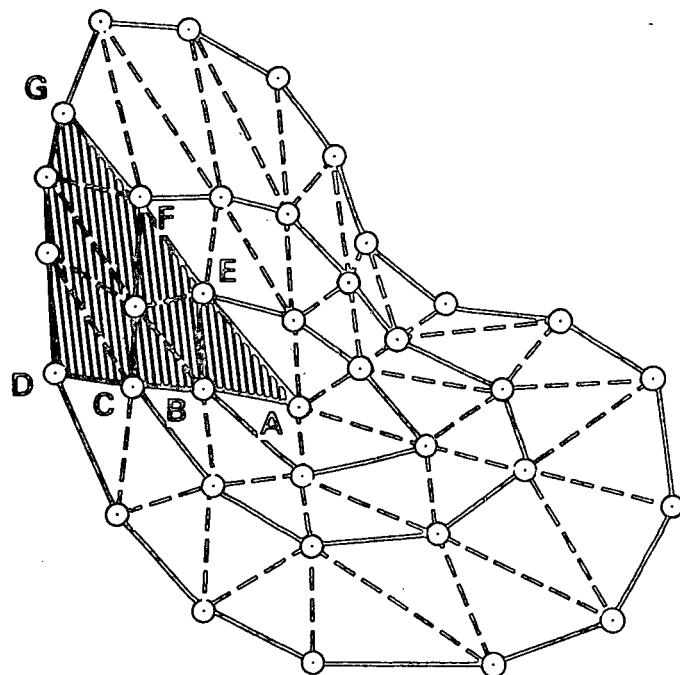


Figure 6. Topology of SNAP break-up.

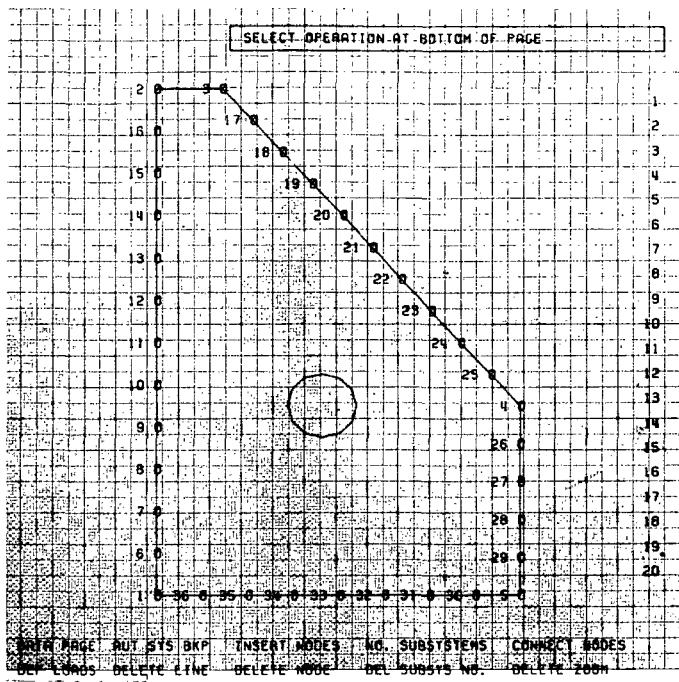


Figure 7. Scope display of boundary.

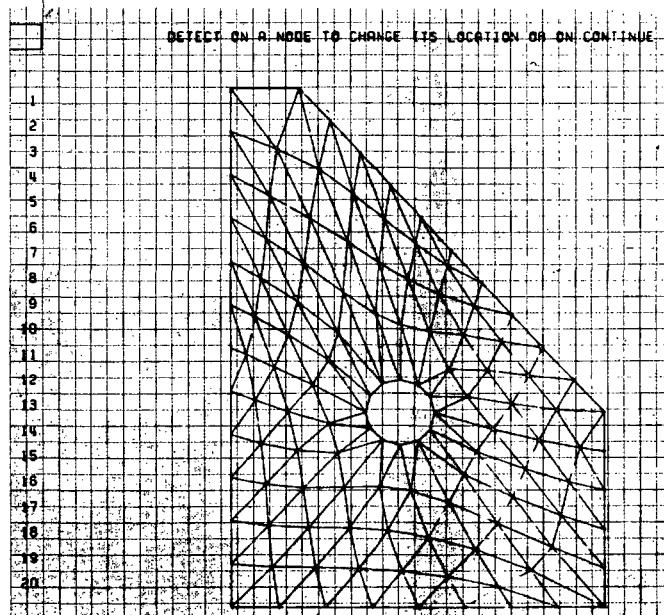


Figure 8. Scope display of coarse SNAP break-up.

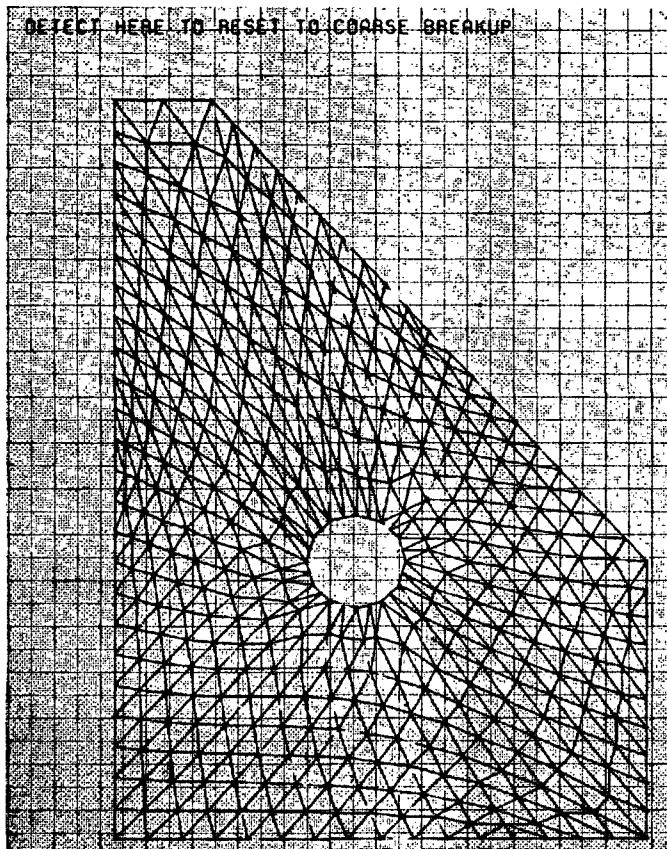


Figure 9. Scope display of fine SNAP break-up.

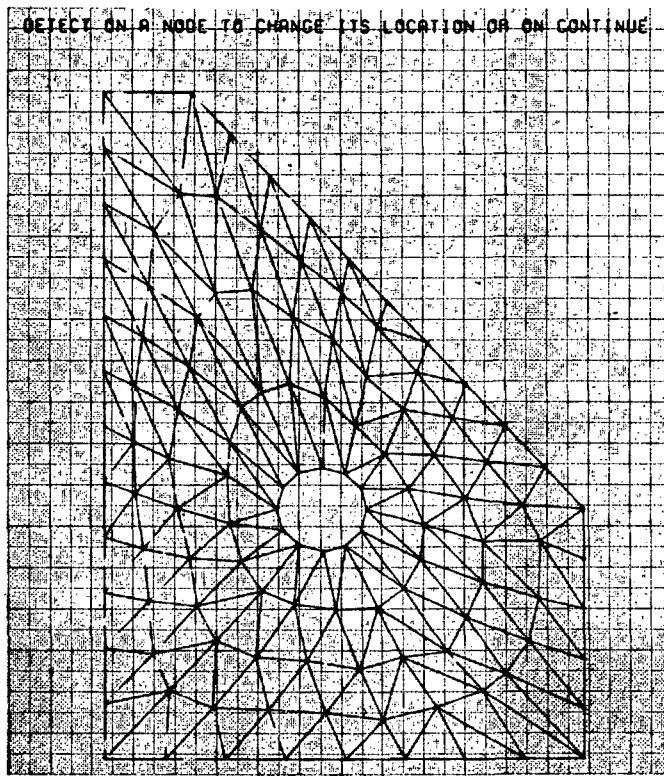


Figure 10. Scope display of adjusted coarse SNAP break-up.

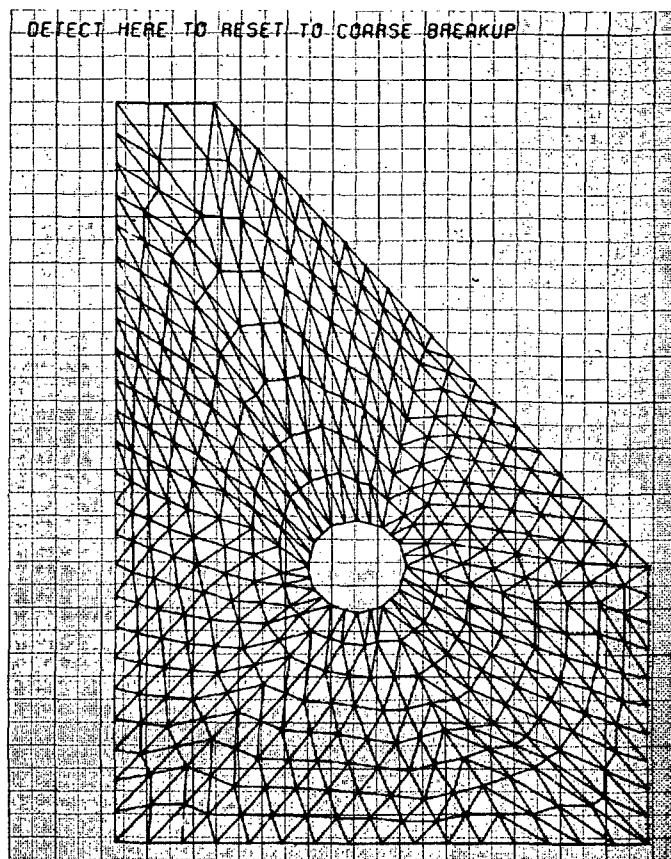


Figure 11. Scope display of adjusted fine SNAP break-up.